



Fig. 3

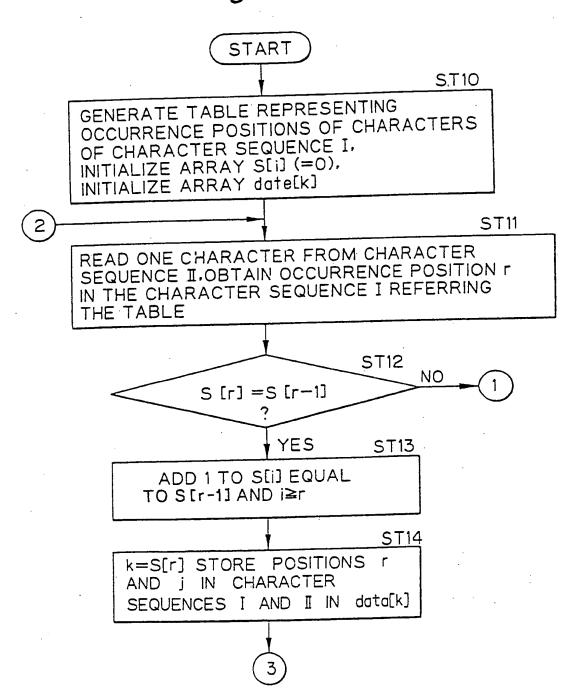




Fig. 4

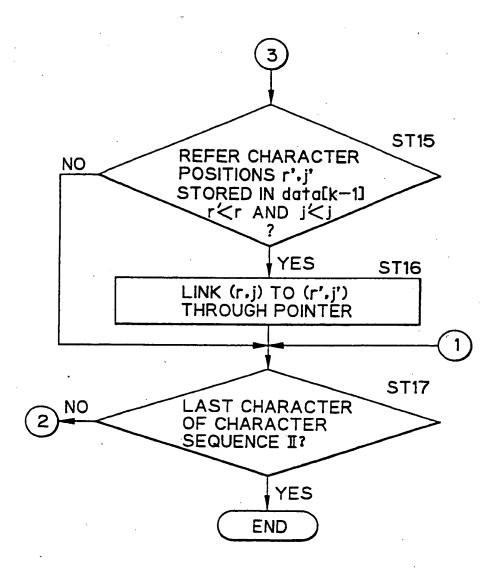
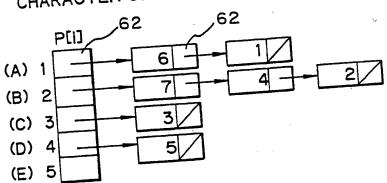


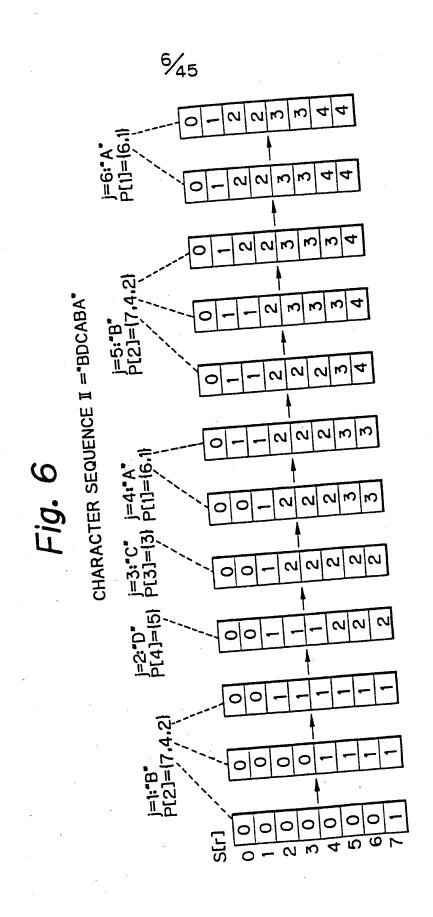


Fig. 5

CHARACTER SEQUENCE I= "ABCBDAB"









7/45

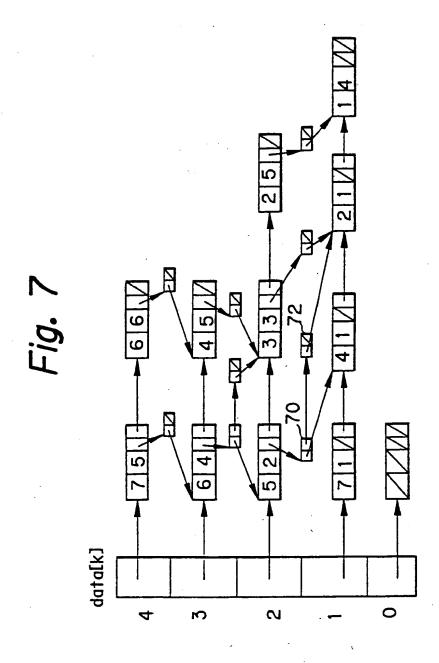




Fig. 8

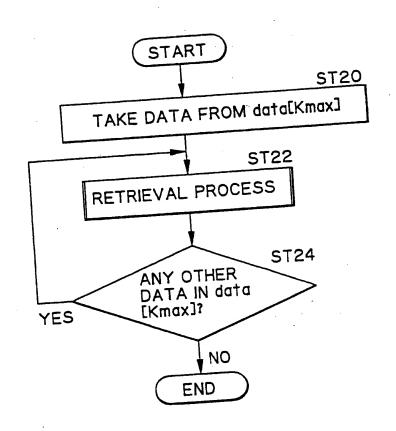




Fig. 9

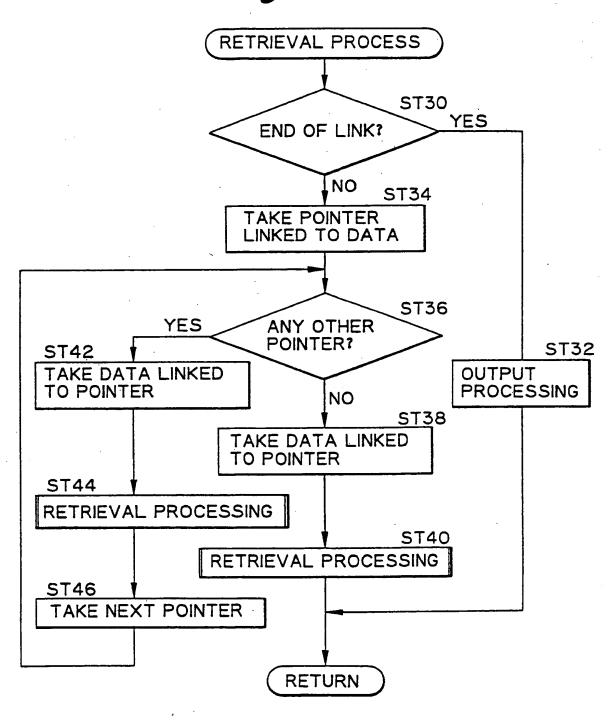




Fig. 10

human bacterium	* GDVEKGKKIFIMKCSQCHTVEGGKHKTGPNLHGLFGRK SEQ ID NO: 1 * EGDAAAGEKVSKKCLACHTFDQGGANKVGPNPNLFGVF SEQ ID NO: 2): 1): 2
SJT	: GD[x3.3]G[x0.1]K[x0.2]K[x4.0]KC[x2.2]CHT[x3.3]GG[x2.2]K GD[x1.4]E[x0.2]K[x0.2]K[x0.4]KC[x2.2]CHT[x3.3]GG[x2.2]K	
homology	: 47%	4



Fig. 11

* MSLAILRVIRLVRVFRIFKLSRHSKGLQILGRTLKASMRELGLLIFFIGVV SEO ID NO: 3 (encinzip, L(6)L(6)L(6)L(6)L Rat



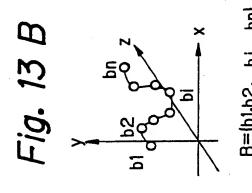
Fig. 12

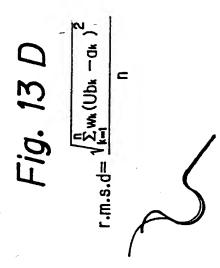
SEQ ID NO: 1 SEQ ID NO: 2

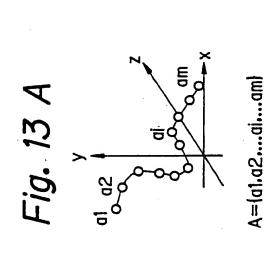
human : GDVEK G K KIFIMKCSQCHTVEKGG KHKTGPNLHGLFGRK ... bacterium : E GDAAAGEKVSK KCLACHTFDQGGANKV GPNPN LFGVF ...

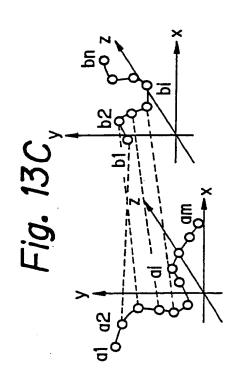














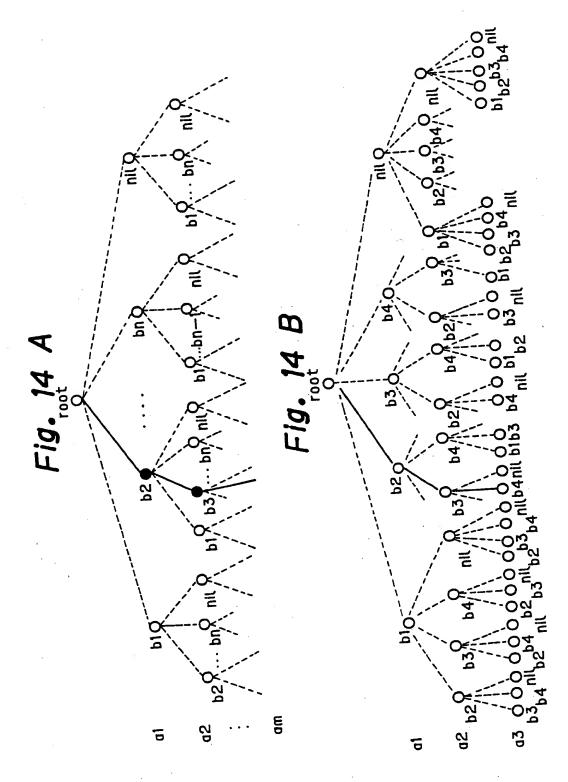
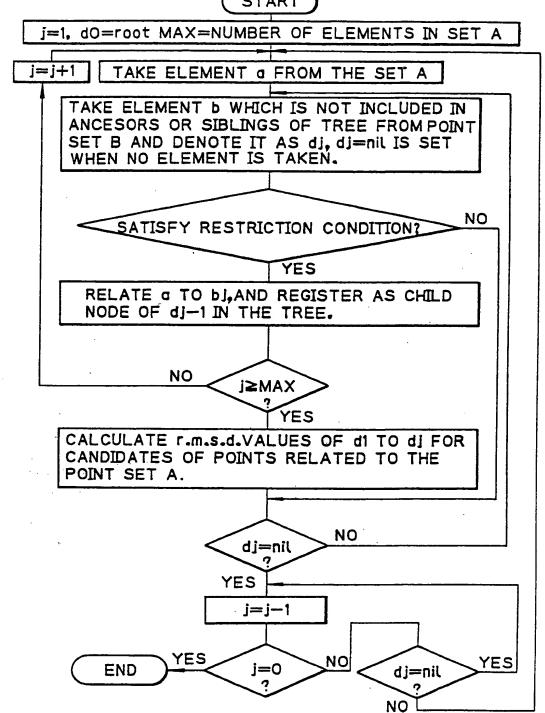




Fig. 15

START



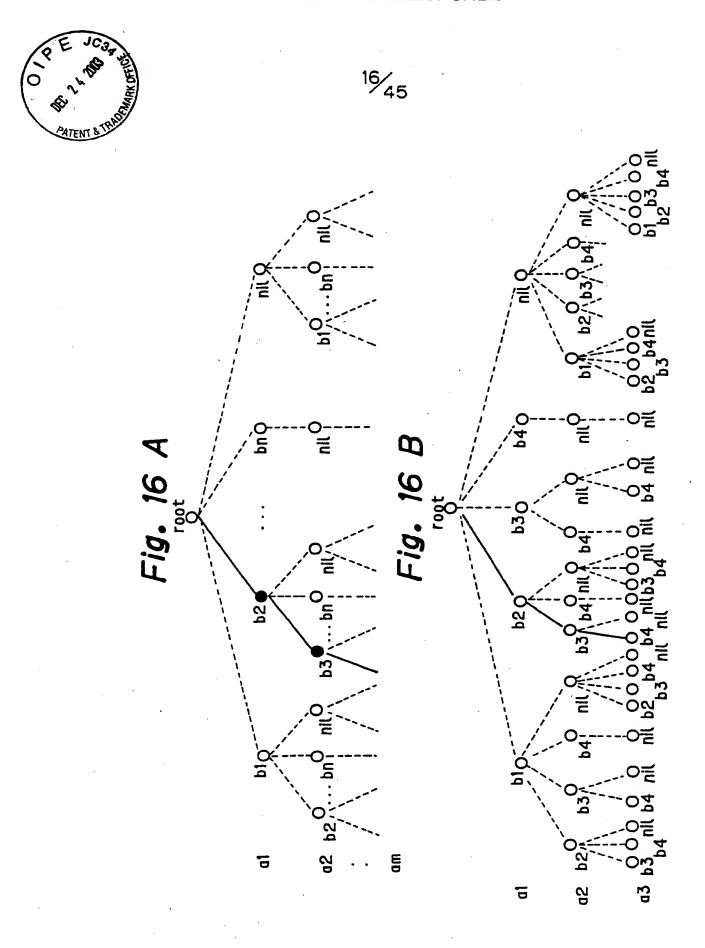
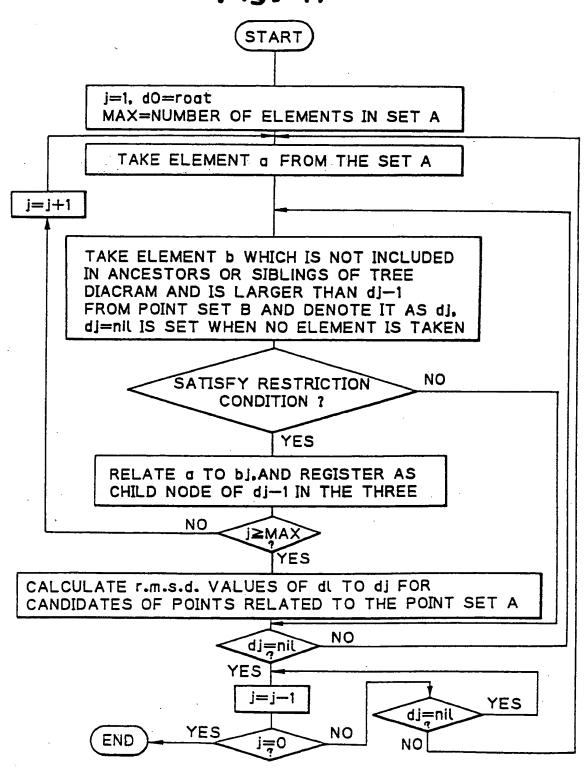




Fig. 17





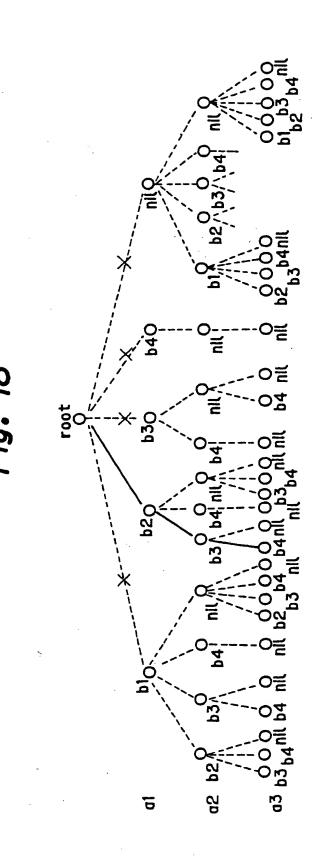
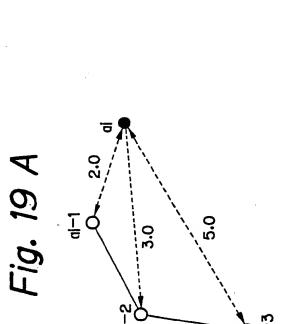
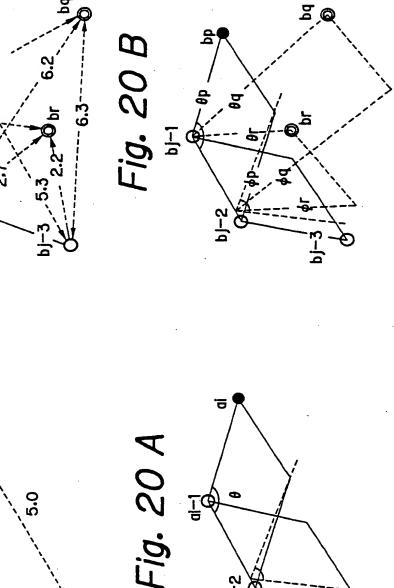






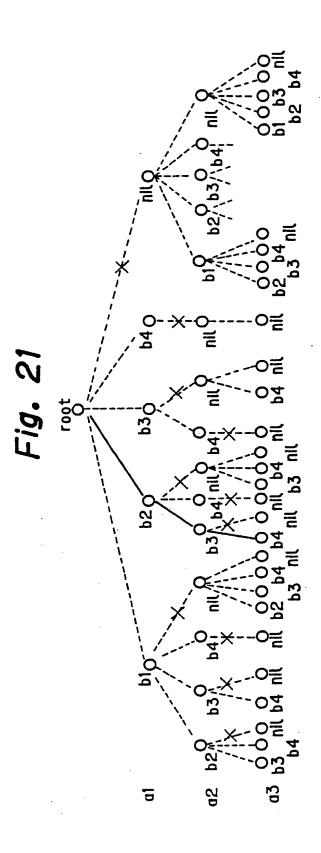
Fig. 19 B





Ē







21/45

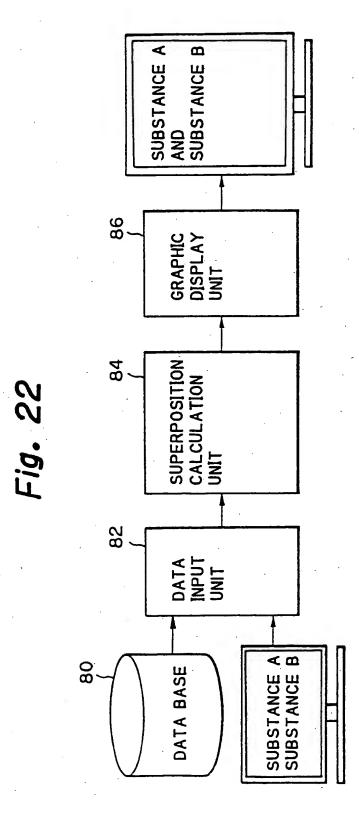




Fig. 23 A

AFSLFDKDGD TEEQIAEFKE VMRSLGQNPT GTITTKELGT VDADGNGTID 21 EAELQDMINE KMKDTDSEEE 41 FPEFLTMMAR DGNGYISAAE IREAFRVFDK KLTDEEVDEM LRHVMTNLGE QVNYEEFVQM 101 IREANIDGDG MTA 141

AMINO ACID SEQUENCE OF CALMODULIN SEQ ID NO: 4

(EXCERPT FROM PDB)

Fig. 23 B

FLSEEMIAEF AMDQQAEARA GGGDISTKEL KAAFDMFDAD PTKEELDAII GTVMRMLGQN IDFEEFLVM EEVDEDGSGT KSEEELADCF VRQMKEDAKG IDIEELGEIL RIFDKNADGF DIEDLMKDSD 101 RATGEHVTEE EFLKMMEGVQ 121 KNNDGRIDFD 141 161

AMINO ACID SEQUENCE OF TROPONIN C SEQ ID NO: 5

(EXCERPT FROM PDB)



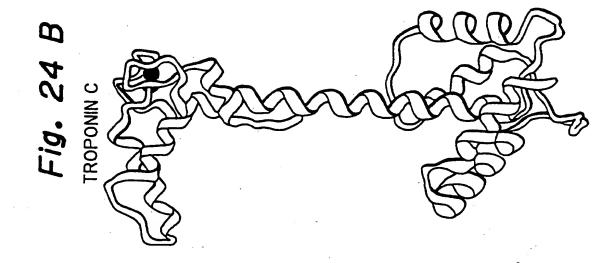
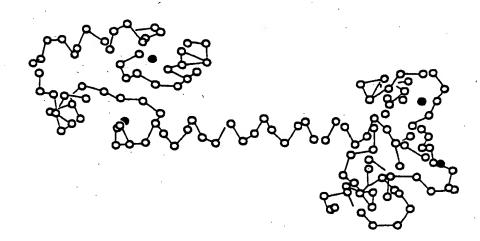


Fig. 24 A CALMODULIN





	O SEQ ID NO: 6	F < target	Y < probe	SEO ID NO: 7	SEO ID NO: 6 < target >	< probe >	SEO ID NO: 7
	100 101 102 103 104 105 106 107 108 109 110	9	g	123	 		
•	108	Q	Z	122	Ø	Z	
	107	A	g	120 121 122 123	<u>~</u>	-	
	901	Z	0	120	_	Σ	
<u> </u>	105	エ	×	6		>	•
81-108 in Calmodulin	104	Q	Q	611 811 211 911	Ш	I	
Sa Imo	103	Ľ.	L_	117	ဗ	<u>~</u>	
<u> </u>	102		>	911	_J	<u></u>	
-108	0	R	K	5	Ш	لنا	
8 ::	00	LL.	ĬĽ.	114	ш	A	7034
	66	ပ	A	113		Ø	567
<u></u>	98		ш	2	0	S	0
Probe site	96 97 98 99	L A	I R	111 112 113			rmsd = 0,567034
·							



Fig. 26

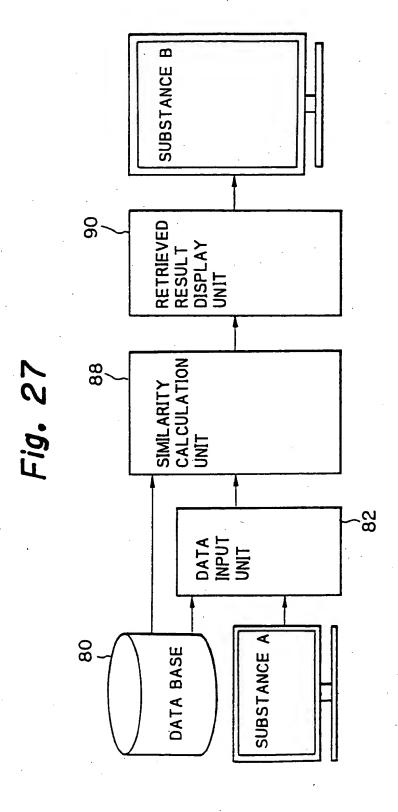
25/45

> SEO ID NO: 8 target > SEQ ID NO: 8 < target > SEQ ID NO. 6 < probe > SEQ ID NO: 7 < probe SEQ ID NO: 9 SEQ ID NO: 8 < target target > probe SEQ ID NO: 7 SEO ID NO: 6 < probe 96 97 98 99 100 101 102 103 104 105 106 107 108 109 110 Probe site = 81-108 and 117-143 in Calmodulin <u>~</u>

rmsd = 0.823665









27/45

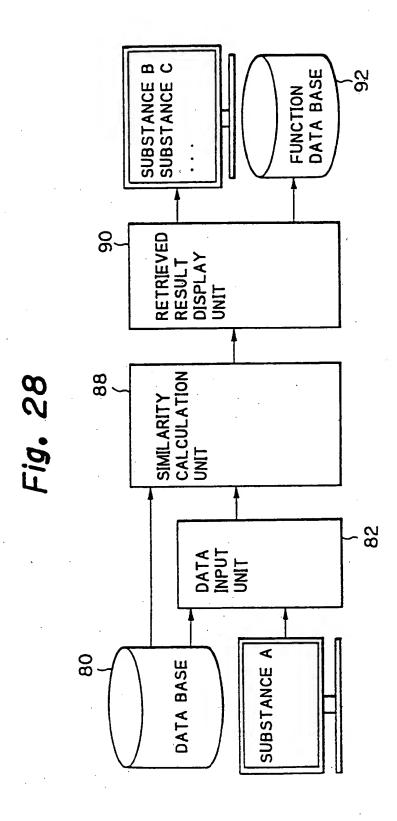


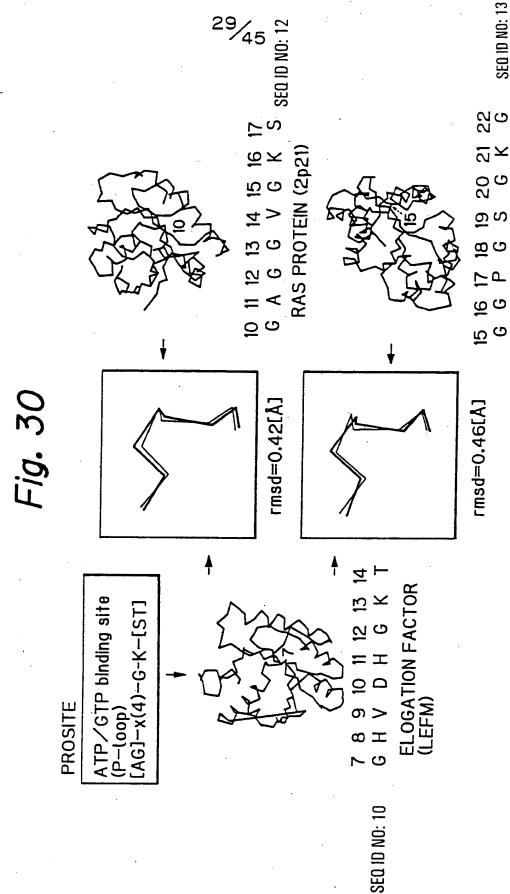


Fig. 29

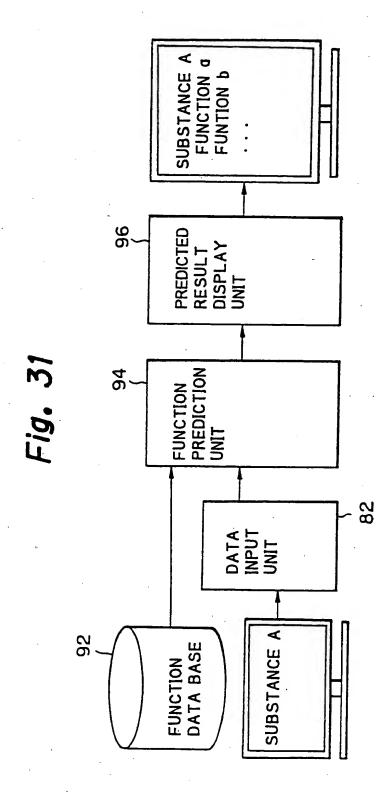
```
SEQ ID NO: 12
SEQ ID NO: 10
                                                                                                                 SEQ ID NO: 11
Seq id no: 10
                                                                   SEQ ID NO: 10
ansansansansansansan ATP/GTP binding site
                                                                  < broke</pre>
                      Probe = (elongation factor)
                                                   ~ 9
```

ADENYLATE KINASE(3ADK)

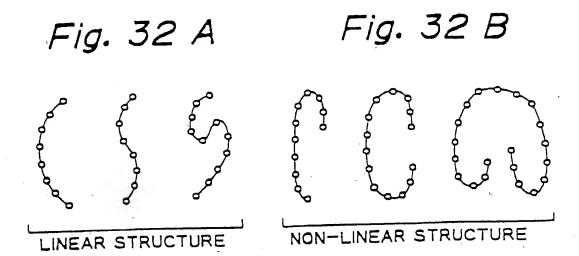












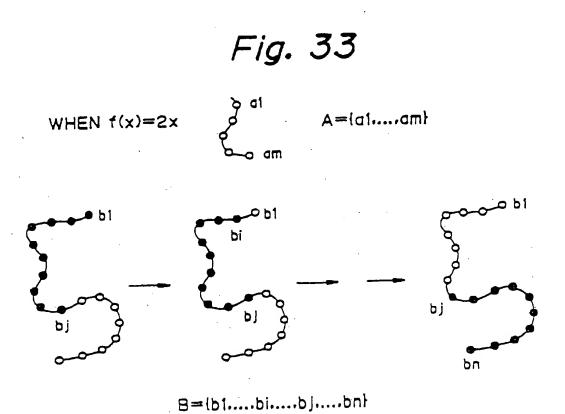




Fig. 34

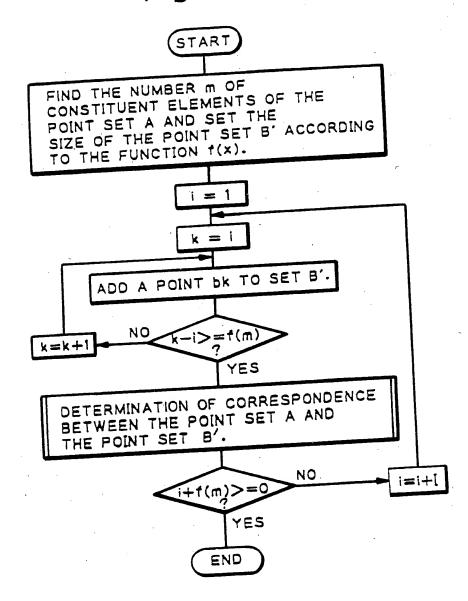
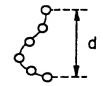


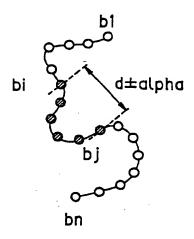


Fig. 35 A



A={a1.a2....am}

Fig. 35 B



B={b1....bi....bj....bn}



Fig. 36

START

PREPARE TABLE OF DISTANCE AMONG POINTS OF THE POINT SETS A,B.

FIND DISTANCE BETWEEN TWO POINTS AT BOTH ENDS OF POINT SET A FROM DISTANCE TABLE AND DENOTE IT AS d.

i = 1

j>i

 $b_j-b_j=d\pm a_j$ p_j m < = j-i < = 2m

SELECT THE ONE HAVING MAXIMUM JOUT OF dj THAT SATISFY THE ABOVE CONDITIONS.

 $B' = \{bi, bi+1, \cdots bj-1, bj\}$

DETERMINATION OF CORRESPONDENCE BETWEEN POINT SET A AND POINT SET B:

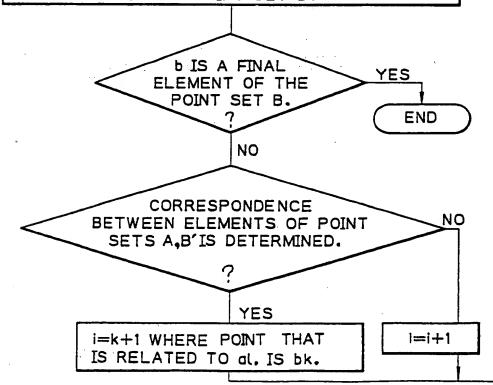




Fig. 37

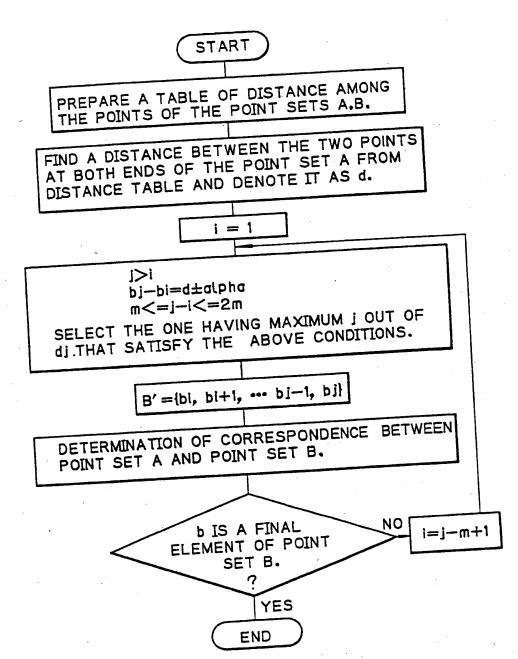




Fig. 38 A

```
IVGGYTCCAN
                 TVPYQVSLNS
   GYHFCGGSLI
                 NSQWVVSAAH
21
   CYKSGIQVRL
                 GEDNI
    EQFISASKS
                 IVHPSYNSNT
    NNDIMLIKL
                 KSAASLNSRV
   ASISLPTSCA
                 SAGTQCLISG
101
    GNTKSSGTS
                 YPDVLKC
   PILSDSSCKS
                 AYPGQITSNM
141
   FCAGYLEGGK
                 DSCQGDSGGP
161
   V V C S G K L Q G I
                 VSWGSGCAQK
181
   NKPGVYTKVC
201
                 NYVSWIKQTI
221
   ASN
             SEQ ID NO: 14
```

AMINO ACID SEQUENCE OF TRYPSIN (EXCERPT FROM PDB)

Fig. 38 B

```
VVGGTEAQRN
                   SWPSQISLQY
   R S G S S W A H T C
M T A A H C V D R E
21
                   GGTLIRQNWV
                   LTFRVVVGEH
61
   NLNQNNGTEQ
                   YVGVQKI
   PYWNTDDVAA
                 GYDIALL
   QSVTLNSYVQ
                   LGVLPRA
101
   LANSPCYITT
121
                   GWGL
                   PTVDYAICSS
141
   SSYWGSTVKN
                  SMVCAGGDGV
161
181
   RSGCQGDSGG
                   PLHCLVNGQY
   AVHGVTSFVS
201
                   RLGCNVTRKP
221
   TVFTRVSAYI
                   SWI
              SEQ ID NO: 15
```

AMINO ACID SEQUENCE OF ELASTASE (EXCERPT FROM PDB)



Fig. 39 A

```
Key site number 36 - 41 in Trypsin

41 42 43 44 45 46

M T A A H C \leftarget \rightarrow SEQ ID NO: 16

V S A A H C \leftarrow probe \rightarrow SEQ ID NO: 17

d = 12.070038 [A]

r.m.s.d. = 0.061077 [A]

The number of atoms in a probe = 6

The number of atoms in PDB = 240

The number of combination = 1

Time = 1sec
```

RETRIEVED RESULTS OF HISTIDINE ACTIVE SITES

Fig. 39 B

```
Key site number 175 - 179 in Trypsin

186 187 188 189 190

G D S G G \left\ target \right\right\ SEQ ID NO: 18

G D S G G \left\ probe \right\right\ SEQ ID NO: 19

d = 8.922721 [A]

r.m.s.d. = 0.092879 [A]

The number of atoms in a probe = 5

The number of atoms in PDB = 240

The number of combination = 1

Time = 1sec
```

RETRIEVED RESULTS OF SERINE ACTIVE SITES



Fig. 40

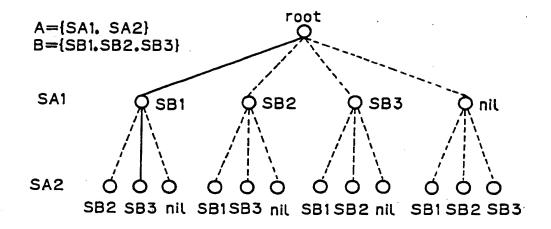
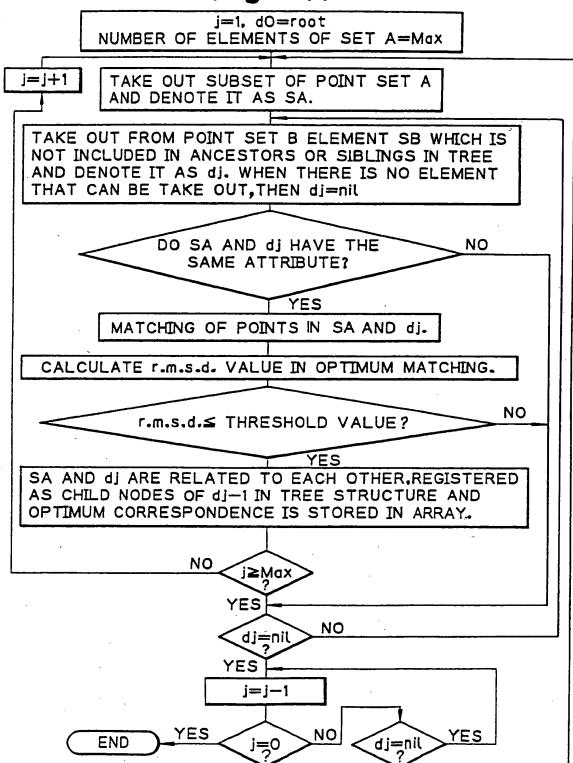




Fig. 41



NO



40/45

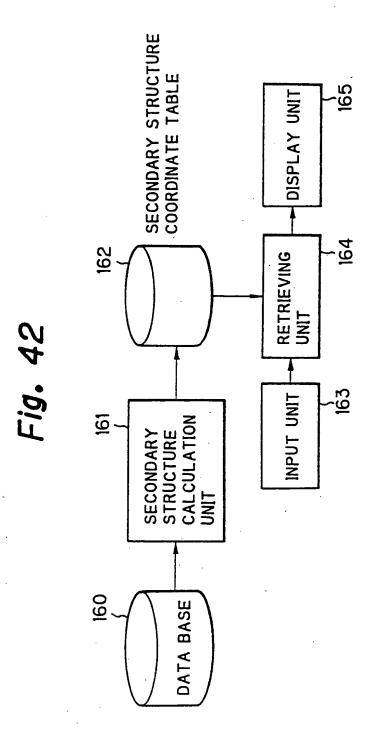
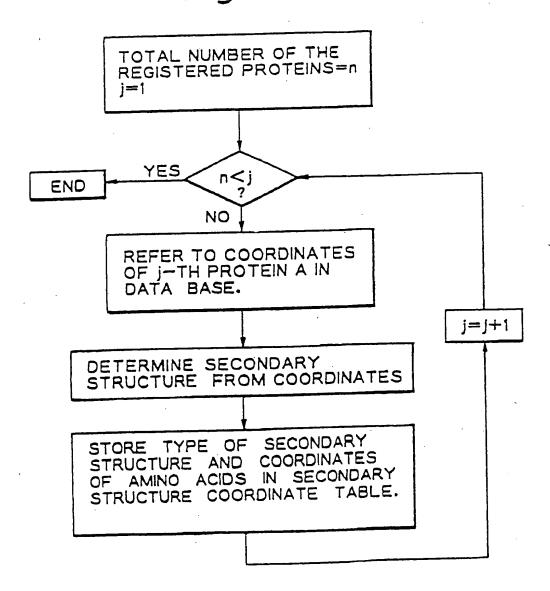




Fig. 43



REPLACEMENT SHEET



42/45

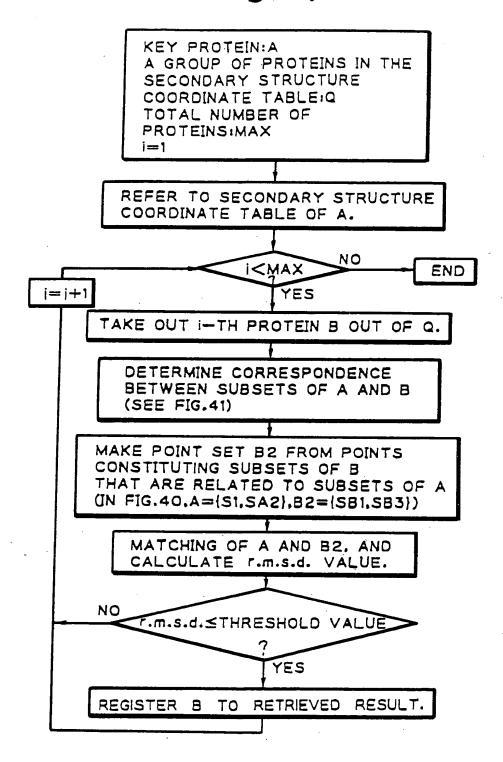
Fig. 44

162

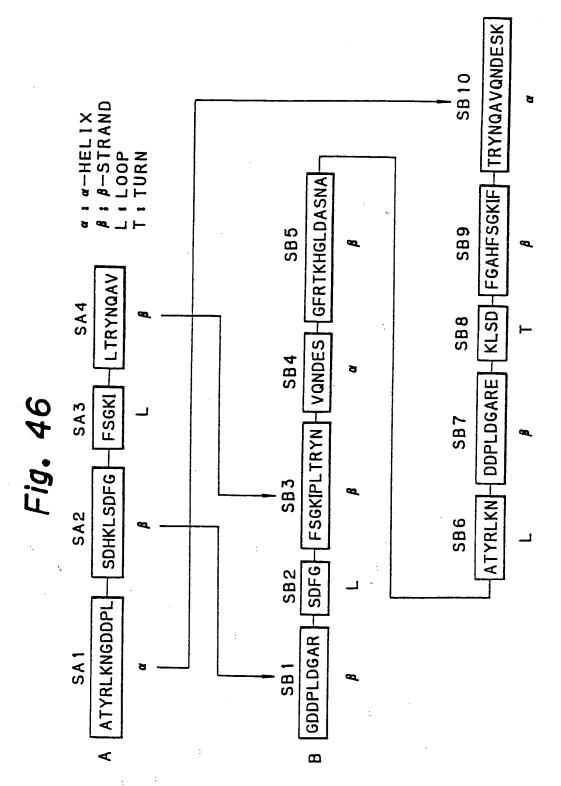
SUBSET	COORDINATES	TYPE
S1	{X1,X2,X3,X4,·····Xa}	α - HELIX
S2	{Xa+1, Xa+2,Xb}	α — HELIX
S3	{Xb+1.Xb+2Xc}	# - SHEET
S4	{Xc+1, Xc+2, Xd}	# — SHEET
Sn	: {X:+1•X:+2•·····Xm}	: 3 — TURN



Fig. 45



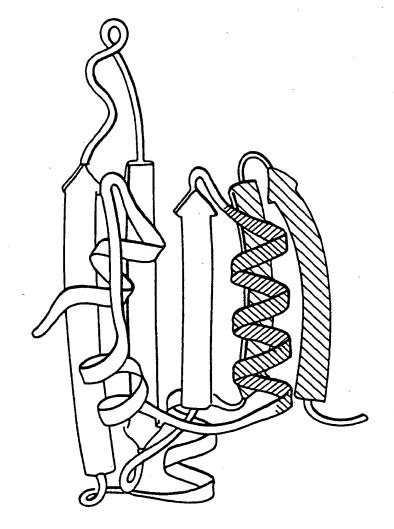




SEQ 10, NO: 20

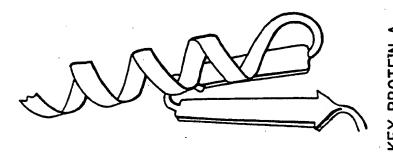






PROTEIN B HAVING A SIMILAR STRUCTURE





KEY PROTEIN A